

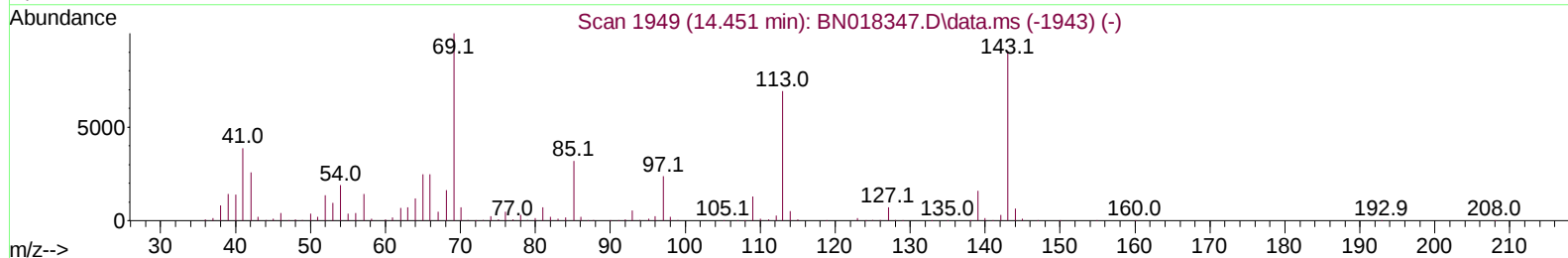
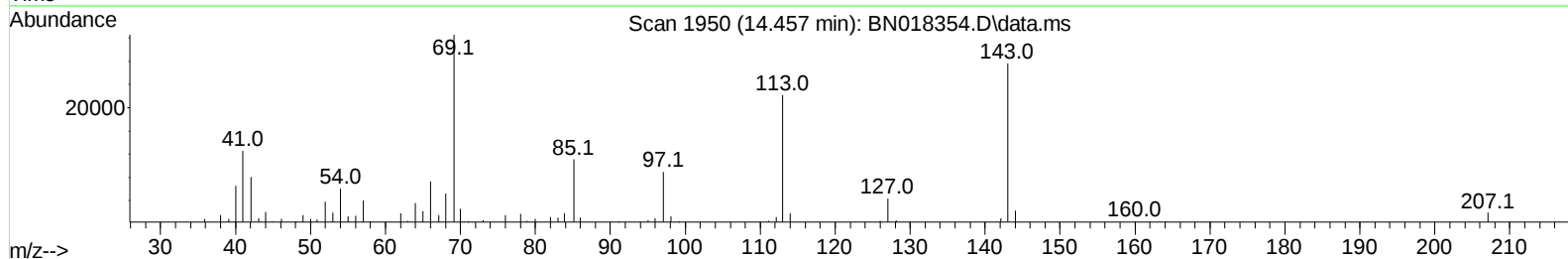
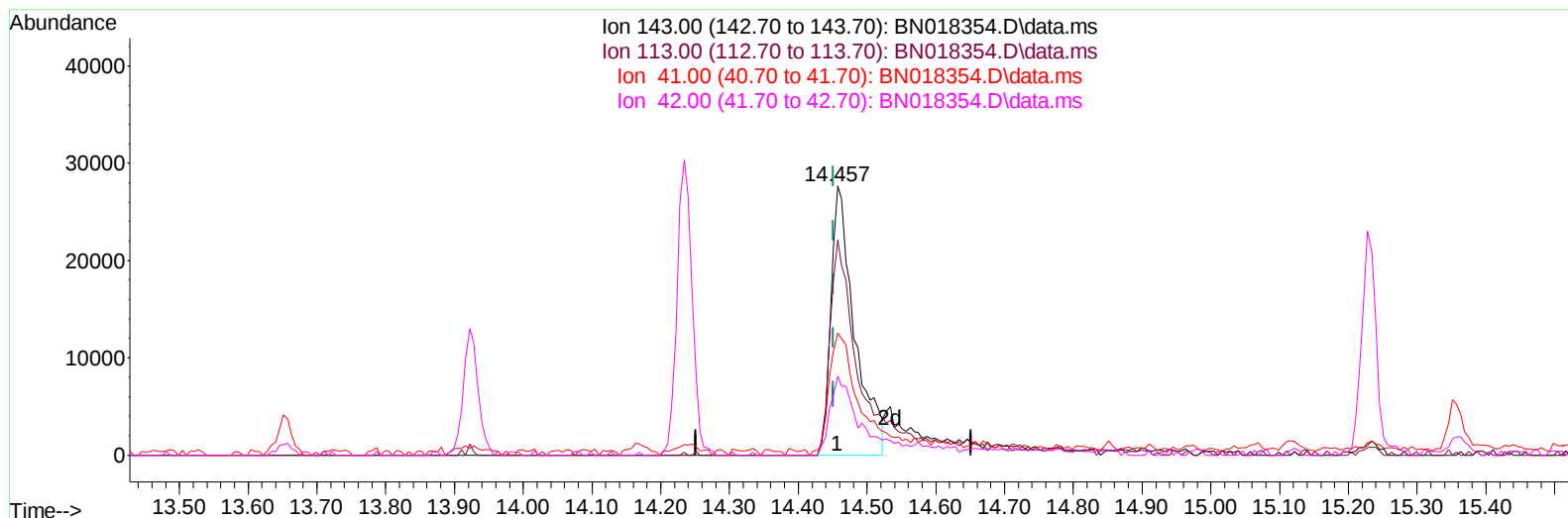
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010622\
Data File : BN018354.D
Acq On : 06 Jan 2022 16:19
Operator : CG/JU
Sample : N1026-02
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGLG6

Manual IntegrationsAPPROVED

Quant Time: Jan 07 09:30:00 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN010622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jan 07 01:31:51 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/07/2022
Supervised By :mohammad ahmed 01/12/2022



TIC: BN018354.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.457min (+ 0.006) 3.82 ng/ul

response 65989

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.00	80.12
41.00	43.80	45.48
42.00	29.00	29.38

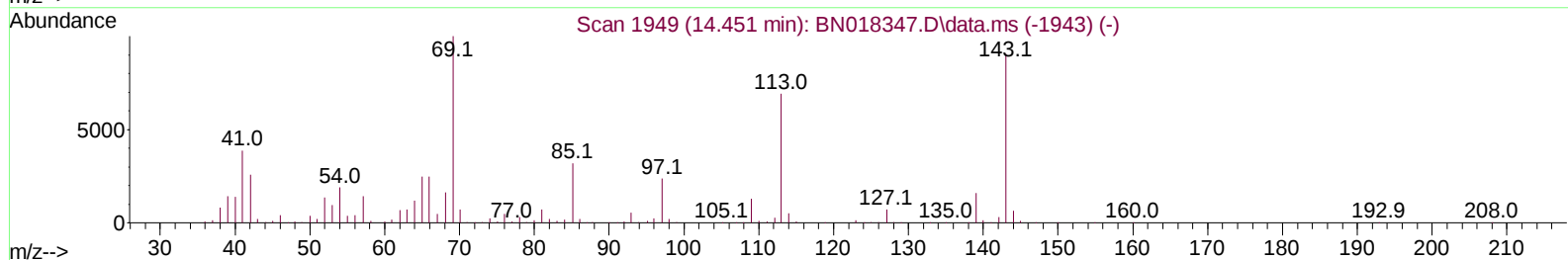
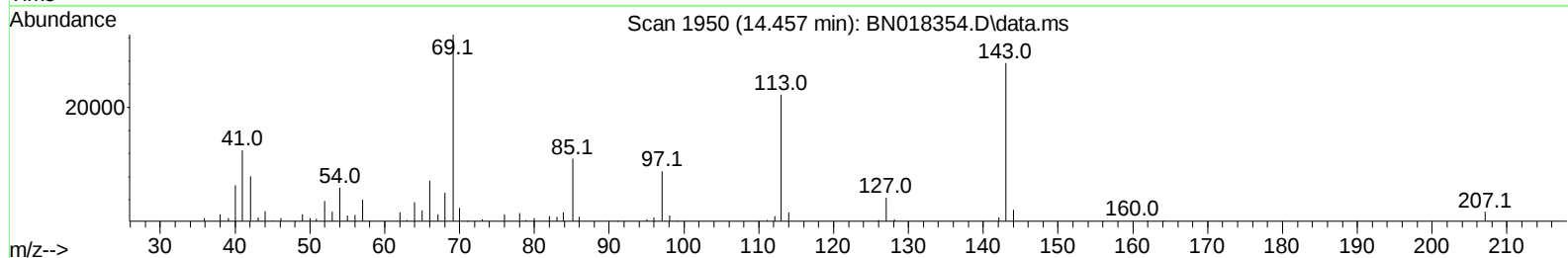
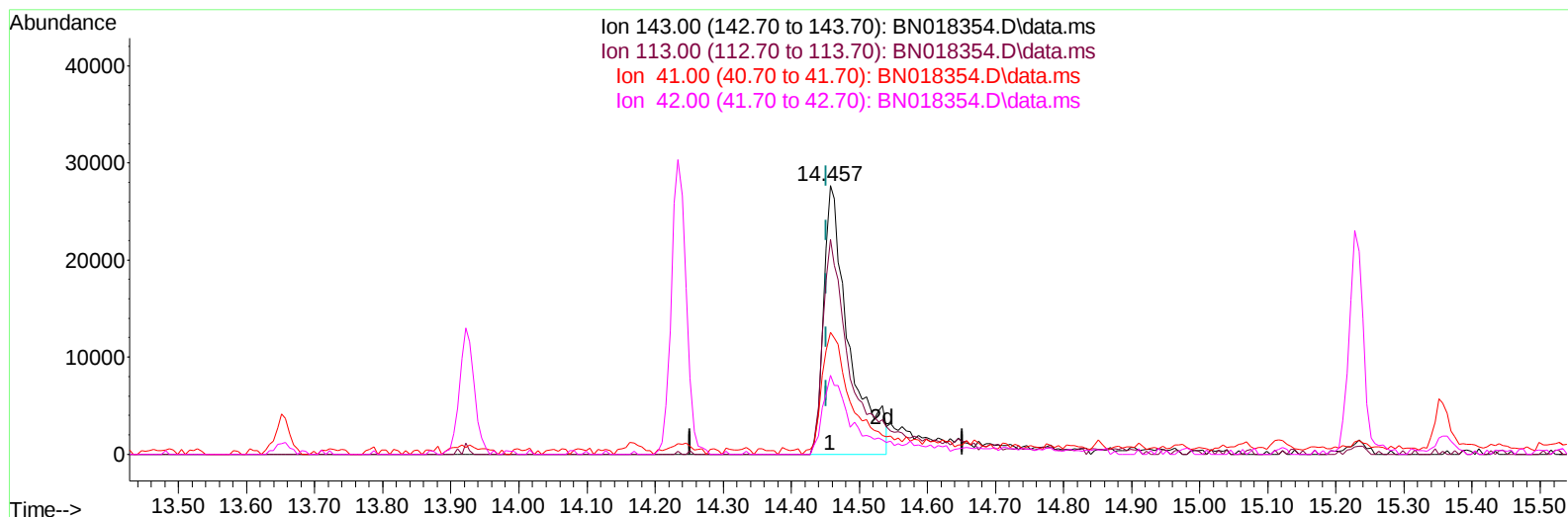
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TIC: BN018354.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.457min (+ 0.006) 4.07 ng/ul m

response 70326

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.00	80.12
41.00	43.80	45.48
42.00	29.00	29.38

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	402399	20.000	ng/ul	0.00
20) Naphthalene-d8	10.369	136	1872631	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.234	164	1245197	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.981	188	2607490	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.175	240	2479819	20.000	ng/ul	# 0.00
88) Perylene-d12	23.398	264	2695898	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.058	96	22402	2.238	ng/uL	0.00
4) Pyridine-d5	3.458	84	747	0.024	ng/ul	0.00
7) Phenol-d5	6.775	99	363324	9.290	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.928	67	305683	12.270	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	306868	10.712	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	302536	9.801	ng/ul	-0.01
21) Nitrobenzene-d5	8.740	128	148917	10.677	ng/ul	0.00
24) 2-Nitrophenol-d4	9.463	143	149923	10.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.005	165	237364	8.763	ng/ul	0.00
31) 4-Chloroaniline-d4	10.510	131	395272	9.495	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	1241515	13.727	ng/ul	0.00
49) Acenaphthylene-d8	13.922	160	1294978	11.326	ng/ul	0.00
54) 4-Nitrophenol-d4	14.457	143	70326m	4.075	ng/ul	0.00
60) Fluorene-d10	15.234	176	1032470	13.530	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	63456	4.301	ng/ul	0.00
73) Anthracene-d10	17.075	188	1676551	13.883	ng/ul	0.00
81) Pyrene-d10	19.375	212	1973342	13.872	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.257	264	1822133	13.581	ng/ul	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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